

```

ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
4-11 8-9 9-10 9-23 11-12 12-13 12-14 19-20 21-22 23-24 23-25 23-26
25-29 26-27 27-28 28-33 29-30 29-31 29-32 33-34 33-35 33-36 37-38
exact bonds :
1-7 7-8 13-15 13-17 14-16 14-18 17-19 18-21 27-37 27-39
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

```

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS
19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS
27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS
35:CLASS 36:CLASS 37:CLASS 38:CLASS 39:CLASS

```

L4 STRUCTURE UPLOADED

```

=> s l4 sss full
FULL SEARCH INITIATED 18:01:07 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 821 TO ITERATE

```

```

100.0% PROCESSED      821 ITERATIONS      0 ANSWERS
SEARCH TIME: 00.00.01

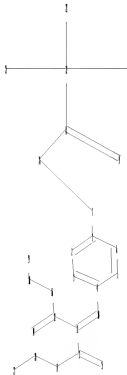
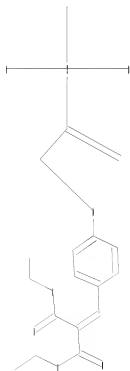
```

L5 0 SEA SSS FUL L4

```

=>
Uploading C:\Documents and Settings\mpepitone\My Documents\ChemDraw\10521629\8.str

```



```

chain nodes :
26
ring nodes :
1 2 3 4 5 6
ring/chain nodes :
7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25
chain bonds :
7-26 8-26
ring/chain bonds :
1-7 4-10 8-22 8-9 10-11 11-12 11-13 12-14 12-16 13-15 13-17 16-18 17-20
18-19 20-21 22-23 22-24 22-25
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
4-10 7-26 8-22 8-9 10-11 11-12 11-13 18-19 20-21 22-23 22-24 22-25
exact bonds :
1-7 8-26 12-14 12-16 13-15 13-17 16-18 17-20
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

```

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS
19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS

```

L6 STRUCTURE UPLOADED

```

=> s l6 sss full
FULL SEARCH INITIATED 18:12:02 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 897 TO ITERATE

```

```

100.0% PROCESSED      897 ITERATIONS      8 ANSWERS
SEARCH TIME: 00.00.01

```

L7 8 SEA SSS FUL L6

```

=> file caplus
COST IN U.S. DOLLARS      SINCE FILE      TOTAL
                        ENTRY      SESSION
FULL ESTIMATED COST      364.54      551.85

```

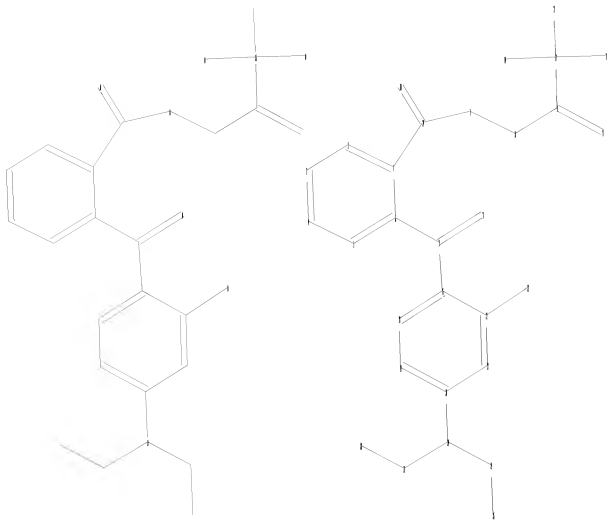
```

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```

chain nodes :
29 30
ring nodes :
1 2 3 4 5 6 16 18 19 20 21 22
ring/chain nodes :
7 8 9 10 11 12 13 14 15 17 23 24 25 26 27 28
chain bonds :
1-29 7-29 29-30
ring/chain bonds :
6-15 7-8 8-9 9-10 9-11 11-12 11-13 11-14 15-16 15-17 20-24 22-23 24-25
24-27 25-26 27-28
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 16-18 16-22 18-19 19-20 20-21 21-22
exact/norm bonds :
6-15 7-29 8-9 9-10 9-11 11-12 11-13 11-14 15-16 25-26 27-28 29-30
exact bonds :

```

1-29 7-8 15-17 20-24 22-23 24-25 24-27
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 16-18 16-22 18-19 19-20 20-21 21-22

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:Atom 17:CLASS 18:Atom
19:Atom 20:Atom 21:Atom 22:Atom 23:CLASS 24:CLASS 25:CLASS 26:CLASS
27:CLASS 28:CLASS 29:CLASS 30:CLASS

L11 STRUCTURE UPLOADED

=> s l11 sss full

FULL SEARCH INITIATED 18:31:08 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 468 TO ITERATE

100.0% PROCESSED 468 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

L12 2 SEA SSS FUL L11

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	358.10	973.63
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-8.00

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=> s l12

L13 1 L12

=> d ibib abs hit str 1-